WHAT IS CLAIMED IS:

A compound of formula I, or a salt or prodrug thereof:

A 1 X W V E - F

 (\mid)

wherein the broken circle represents two non-adjacent double bonds in any position in the five-membered ring;

two, three or four of V, W, X, Y and Z represent nitrogen and the remainder represent carbon provided that, when two of V, W, X, Y and Z represent nitrogen and the remainder represent carbon, then the said nitrogen atoms are in non-adjacent positions within the five-membered ring;

A¹ is selected from the group consisting of, hydrogen, hydrocarbon, a heterocyclic group, halogen, cyano, trifluoromethyl, -OR*, -SR*, -NR*R*, -NR*COR*, -NR*CO₂R*, -NR*SO₂R*, and -NR²CTNR*R*;

A² represents a non-bonded electron pair when four of V, W, X, Y and Z represent nitrogen and the other represents carbon; or, when two or three of V, W, X, Y and Z represent nitrogen and the remainder represent carbon, A² is selected from the group-consisting of hydrogen, hydrocarbon, a heterocyclic group, halogen, cyano, trifluoromethyl, -OR^x, -SR^x, -NR^xR^y, -NR^xCOR^y, -NR^xCO₂R^y, -NR^xSO₂R^y, and -NR^zCTNR^xR^y;

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E represents a bond or a straight or branched alkylene chain containing from 1 to 4/carbon atoms;

F represents a group of formula

U represents nitrogen or C-R²;

B represents oxygen, sulphur or N-R3;

R¹ represents -CH₂.CHR⁴.NR⁶R⁷ or a group of

15 formula

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$$N-R^5$$
, $N-R^5$

in which the broken line represents an optional chemical bond;

 R^2 , R^3 , R^4 , R^5 , R^6 and R^7 independently represent hydrogen or C_1 , alkyl;

 R^{x} and R^{y} independently represent hydrogen, hydrocarbon or a heterocyclic group, or R^{x} and R^{y} together represent a φ_{2-6} alkylene group;

R^z/represents hydrogen, hydrocarbon or a heterocyclic group;

Trepresents oxygen, sulphur or a group of formula =N/G; and

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G represents hydrocarbon, (a/heterocyclic group or an electron-withdrawing group.

A compound according to claim 1 5 represented by formula IIA, and salts and prodrugs thereof:

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X1 represents nitrogen or A12-C; n is zero, 1/2 or 3; B¹ represents oxygen, sulphur or N-R¹³;

A¹¹ and A¹² are independently selected from the 20 group consisting of c_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-7} cycloalkyl, aryl, aryl(C_{1-6}) alkyl, C_{3-7} heterocycloalkyl, heteroaryl and heteroaryl(C1-6)alkyl, any of which group's may be optionally substituted; and hydrogen, halogen, cyano, trifluoromethyl, C1-6 alkoxy, C₁₋₆ alkylthio and -NR^xR^y; 25

 R^{12} , R^{13} , R^{14} , R^{16} and R^{17} independently represent hydrogen or $C_{1-6}/alkyl$; and

Rx and Ry independently represent hydrogen, hydrocarbon or a heterocyclic group, or Rx and Ry together represent a C₂₋₆ alkylene group.

A compound according to claim 1 represented/by formula IIB, and salts and prodrugs thereof:

$$A^{21} \longrightarrow_{Y^1 = N} (CH_2)_n$$

$$= NR^{26}R^{27}$$

$$= R^{24}$$

$$= R^{24}$$

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Y¹ represents nitrogen or A²²-C;

n is zero, 1, 2 or 3;

B² represents oxygen, sulphur or N-R²³;

A²¹ and A²² are independently selected from the group consisting of C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₇ cycloalkyl, aryl, aryl(C₁₋₆)alkyl, C₃₋₇ heterocycloalkyl, heteroaryl and heteroaryl(C₁₋₆)alkyl, any of which groups may be optionally substituted; and hydrogen, halogen, cyano, trifluoromethyl, C₁₋₆ alkoxy,

20 C₁₋₆ alkylthio and -MR^xR^y;

 R^{22} , R^{23} , R^{24} , R^{26} and R^{27} independently represent hydrogen or C_{1-6} alkyl; and

 R^x and R^y independently represent hydrogen, hydrocarbon or a heterocyclic group, or R^x and R^y together represent a \mathcal{C}_{2-6} alkylene group.

/4. A compound according to claim 1 represented by formula IIC, and salts and prodrugs thereof:

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$$\begin{array}{c|c}
 & A & 3 & 1 \\
 & N & (CH_2)_n & R^{31} \\
 & \downarrow 2 & = 1 \\
 & \downarrow 2 & = 1
\end{array}$$
(1)(C)

wherein

Y² represents nitrogen or A³²-C;

Z¹ represents nitrogen or CH;

n is zero, 1, ½ or 3;

B³ represents/oxygen, sulphur or N-R³³;

A³¹ and A³² are independently selected from the group consisting of C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₇ cycloalkyl, aryl, aryl(C₁₋₆)alkyl, C₃₋₇ heterocycloalkyl, heteroaryl and heteroaryl(C₁₋₆)alkyl, any of which groups may be optionally substituted; and hydrogen, halogen, cyano, trifluoromethyl, C₁₋₆ alkoxy, C₁₋₆ alkylthio and -NR^xR^y;

R³¹ represents -CH₂.CHR³⁴.NR³⁶R³⁷ or a group of formula

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$$N-R^{35}$$
 or $N-R^{35}$;

 R^{32} , R^{33} , R^{34} , R^{35} , R^{36} and R^{37} independently represent hydrogen or C_{1-6} alkyl; and

 R^x and R^y independently represent hydrogen, hydrocarbon or a heterocyclic group, or R^x and R^y together represent a C_{2-6} alkylene group.

5. A compound according to claim 1 represented by formula IID, and salts and prodrugs thereof:

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$$\begin{array}{c|c}
A^{41} & W^{1} \\
N = N
\end{array}$$

$$\begin{array}{c|c}
R^{41} \\
R^{42}
\end{array}$$

$$\begin{array}{c|c}
R^{42}
\end{array}$$

wherein

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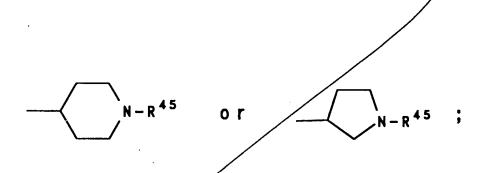
W¹ represents nitrogen or C-A⁴²; n is zero, 1, 2 or 3;

B4 represents oxygen, sulphur or N-R43;

A⁴¹ and A⁴² are independently selected from the group consisting of C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₇ cycloalkyl, aryl, aryl(C₁₋₆)alkyl, C₃₋₇ heterocycloalkyl, heteroaryl and heteroaryl(C₁₋₆)alkyl, any of which groups may be optionally substituted; and hydrogen, halogen, cyano, trifluoromethyl, C₁₋₆ alkoxy, C₁₋₆ alkylthio and -NR^xR^y;

R⁴¹ represents -CH₂.CHR⁴⁴.NR⁴⁶R⁴⁷ or a group of formula

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R⁴², R⁴³, R⁴⁴, R⁴⁵, R⁴⁶ and R⁴⁷ independently

represent hydrogen or C₁₋₆ alkyl; and

R^x and R^y independently represent hydrogen,

hydrocarbon or a heterocyclic group, or R^x and R^y together

represent a C₂₋₆ alkylene group.

6. A compound according to claim 1 selected from:

2-[5-(2-benzyltetrazol-5-ylmethyl)-1H-indol-3-yl]ethylamine;

2-[5-(1-benzyltetrazol-5-ylmethyl)-1H-indol-3-yl]ethylamine;

N, N-dimethyl-2-[5-(1-methyltetrazo1-5-ylmethyl)-1H-indol-3-yl]ethylamine;
N, N-dimethyl-2-[5-(2-methyltetrazo1-5-ylmethyl)-1H-indol-3-yl]ethylamine;

N,N-dimethyl-2-[5-(1,2,4-triazol-1-ylmethyl)-1H-indol-3-yl]ethylamine;

N, N-dimethyl-2-[5-(tetrazol-2-ylmethyl)-1H-indol-3-yl]ethylamine;

N, N-dimethyl-2-[5-(tetrazol-1-ylmethyl)-1H-indol-3-

yl]ethylamine;
N,N-dimethyl-2-[5-(1-methyl-1,2,4-triazol-5-ylmethyl)-1Hindol-3-yl]ethylamine;
N,N-dimethyl-2-[5-(1-methyl-1,2,4-triazol-3-ylmethyl)-1H-

indol-3-yl]ethylamine;

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N, N-dimethyl-2-[5-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-triazol-1-ylmethyl)-1H-ind\phil-3-(1,2,3-tria
               yl]ethylamine;
               3-(2-aminoethyl)-5-(1-methyltetrazol-5-yl)-
               benzo[b]thiophene;
                3-(2-aminoethyl)-5-(2-methyltetrazol-5-yl)-
               benzo[b]thiophene;
                3-[2-(N,N-dimethylamino)ethyl]-5-(2-methylt/etrazol-5-
               yl)benzo[b]thiophene;
                N, N-dimethyl-2-[5-(2-methylimidazol-1-y/methyl)-1H-indol-
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                3-yl]ethylamine;
                N, N-dimethyl-2-[5-(imidazol-1-ylmethyl)-1H-indol-3-
                yl]ethylamine;
                N, N-dimethyl-2-[5-(2-methylimidaz\phil-1-yl)-1H-indol-3-
               yl]ethylamine;
               N, N-dimethyl-2-[5-(2-ethyltetrazol-5-ylmethyl)-1H-indol-
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                3-yl]ethylamine;
                N, N-dimethyl-2-[5-(1-ethyltetrazol-5-ylmethyl)-1H-indol-
                3-yl]ethylamine;
               N, N-dimethyl-2-[5-(1,2,47triazol-1-yl)-1H-indol-3-
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                yl]ethylamine;
         - 1-methyl-4-[5-(2-methylimidazol-1-yl)-1H-indol-3-
                yl]piperidine;
                1-methyl-4-[5-(1,2,4-triazol-1-ylmethyl)-1H-indol-3-
                yl]piperidine;
25 — 4-[5-(2-methylimidazol-1-yl)-1H-indol-3-yl]piperidine;
         ->4-[5-(1,2,4-triazol-1-ylmethyl)-1H-indol-3-yl]piperidine;
                 3-[5-(2-methyl/imidazol-1-yl)-1H-indol-3-yl]pyrrolidine;
                1-\text{methyl}-3-[5/-(2-\text{methylimidazol}-1-\text{yl})-1\text{H-indol}-3-
                yl]pyrrolidine;
                4-[5-(imidazol-1-yl)-1H-indol-3-yl]piperidine;
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                4-[5-(1,2/,3-triazol-1-yl)-1H-indol-3-yl]piperidine;
                 1-methy1/-4-[5-(imidazol-1-yl)-1H-indol-3-yl]piperidine;
                 1-\text{methyl}-4-[5-(1,2,3-\text{triazol}-1-\text{yl})-1\text{H-indol}-3-
                yl]piperidine;
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1-methyl-3-[5-(1,2,3-triazol-1-yl)-1H-indol-3yl]pyrrolidine; 1-methyl-3-[5-(2-methylimidazol-1-ylmethyl)-1H-indol-3yllpyrrolidine; 1-methyl-3-[5-(imidazol-1-yl)-1H-indol-3-yl]pyrrolidine; 5 1-methyl-3-[5-(1,2,4-triazol-1-ylmethyl)-1H-indol-3yl]pyrrolidine; 1-methyl-3-[5-(imidazol-1-y/methy/)-1H-indol-3yl|pyrrolidine; N, N-dimethyl-2-[5-(2-aminoimidazol-1-yl)-1H-indol-3-10 yl]ethylamine; N, N-dimethyl-2-[5-(2-aminoimidazol-1-ylmethyl)-1H-indol-3-yl]ethylamine; N-methyl-2-[5-(1,2,4-ty/iazol-1-ylmethyl)-1H-indol-3-15 yllethylamine;

7. A pharmaceutical composition comprising an effective amount of a compound according to claim 1 in association with a pharmaceutically acceptable carrier or excipient.

8. A method for the treatment and/or prevention of clinical conditions for which a selective agonist of 5-HT₁-like receptors is indicated, which method comprises administering to a patient in need of such treatment an effective amount of a compound according to claim 1.

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and salts and prodrugs thereof.